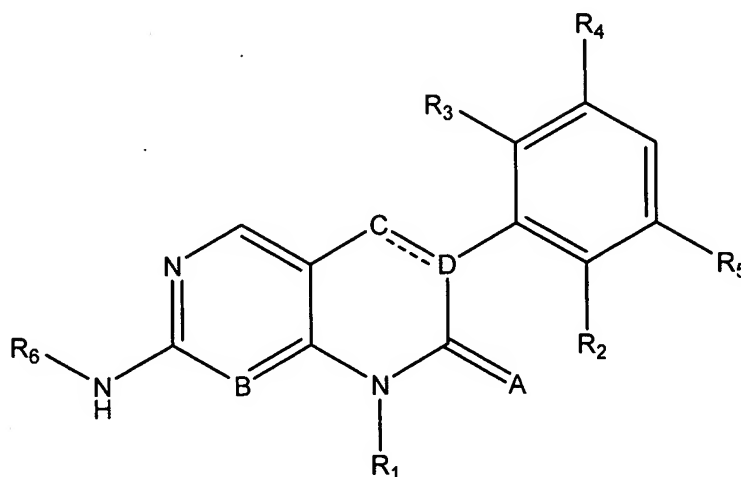


**Amendments to the Claims:**

This listing of claims will replace all prior versions and listings of claims in the application:

**Listing of Claims:**

1. (Currently Amended) A compound of the formula:



wherein:

A is O,  $\text{NH}_2$ ,  $\text{NH}(\text{C}_1\text{-C}_6 \text{ alkyl})$ ,  $\text{N}(\text{C}_1\text{-C}_6 \text{ alkyl})_2$ , or  $\text{-NHC(O)NHR}_{12}$ ;

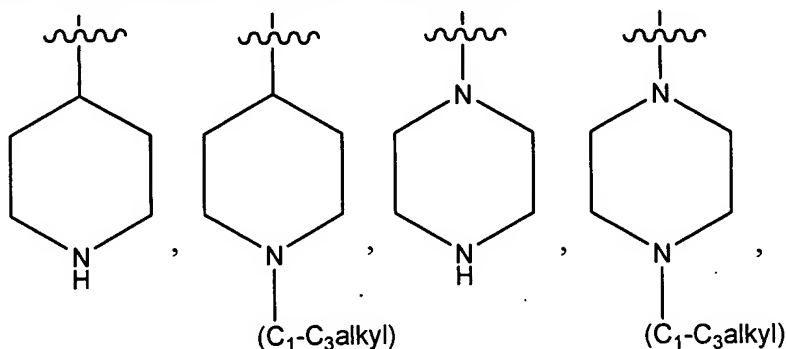
$\text{R}_{12}$  is  $\text{C}_1\text{-C}_6$  straight or branched chain alkyl, or  $\text{-(CH}_2\text{)}_n\text{-C}_3\text{-C}_8$  cycloalkyl ring; n is an integer of from 1 to 3;

~~B, C and D are independently selected from CH or N, with the proviso that C and D are not both N; B is N, C is CH, D is N.~~

$\text{R}_1$  is selected from the group of  $\text{C}_1\text{-C}_6$  straight or branched chain alkyl, optionally substituted by  $\text{-COOH}$ , or;

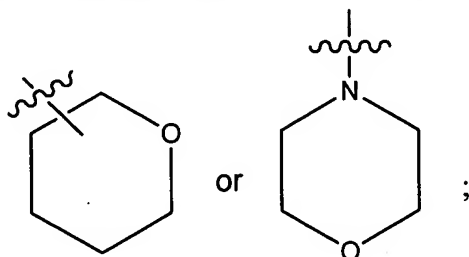
a) a phenyl, benzyl or  $\text{C}_3\text{-C}_8$  cycloalkyl ring, or  $\text{-CH}_2\text{-C}_3\text{-C}_8$  cycloalkyl ring, with the phenyl, benzyl or cycloalkyl rings being optionally substituted by 1 or 2  $\text{COOH}$  or  $\text{-CH}_2\text{-COOH}$  groups; or

b) a piperidine or piperazine moiety selected from group of:



the rings of the piperidine or piperazine moieties being optionally substituted by 1 or 2 COOH or -CH<sub>2</sub>-COOH groups; or

c) a tetrahydropyran or morpholine moiety of the formulae:



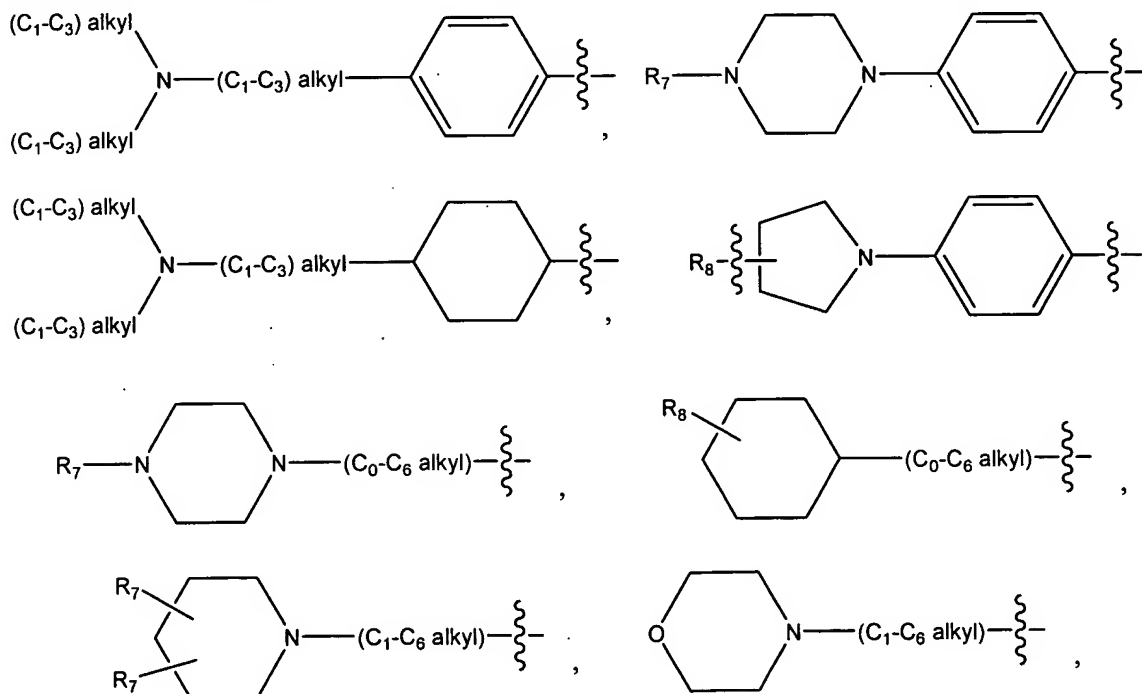
R<sub>2</sub> is H, Cl or F;

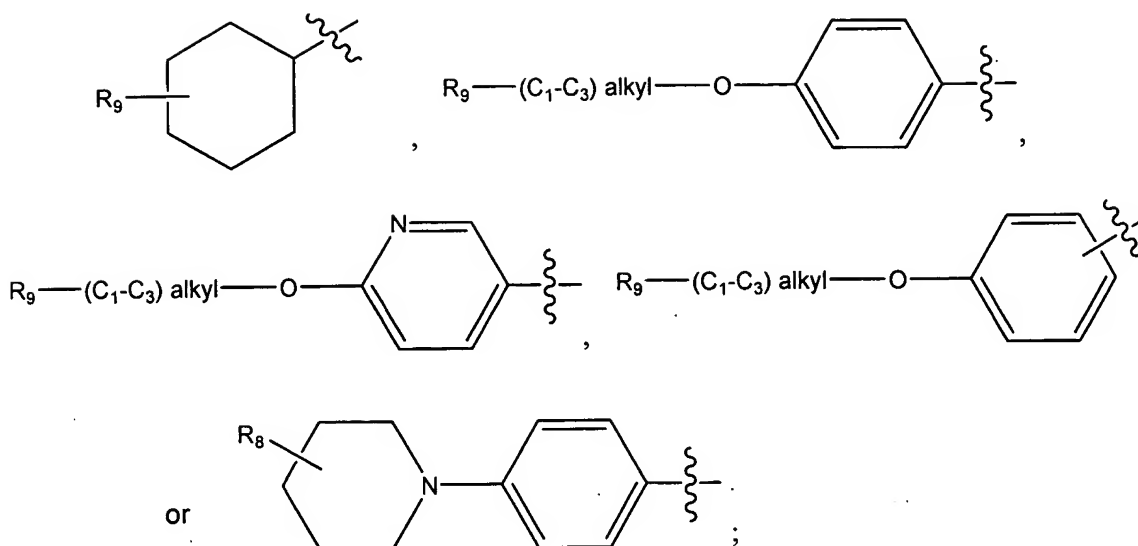
R<sub>3</sub> is H, Cl or F, with the proviso that at least one of R<sub>2</sub> or R<sub>3</sub> is F;

R<sub>4</sub> is H, OH, -OCH<sub>3</sub>, or -OCH<sub>2</sub>CH<sub>3</sub>, with the proviso that, if R<sub>4</sub> is H, R<sub>2</sub> and R<sub>3</sub> are not H;

R<sub>5</sub> is -OCH<sub>3</sub>, or -OCH<sub>2</sub>CH<sub>3</sub>;

R<sub>6</sub> is selected from the group of H, -(C<sub>1</sub>-C<sub>5</sub> alkyl)-NH<sub>2</sub>, -(C<sub>1</sub>-C<sub>5</sub> alkyl)-NH-(C<sub>1</sub>-C<sub>3</sub> alkyl)-R<sub>11</sub>, -(C<sub>1</sub>-C<sub>5</sub> alkyl)-N-(C<sub>1</sub>-C<sub>3</sub> alkyl-R<sub>11</sub>)<sub>2</sub>, -O-(C<sub>1</sub>-C<sub>5</sub> alkyl)-NH<sub>2</sub>, -O-(C<sub>1</sub>-C<sub>5</sub> alkyl)-NH-(C<sub>1</sub>-C<sub>3</sub> alkyl)-R<sub>11</sub>, -O-(C<sub>1</sub>-C<sub>5</sub> alkyl)-N-(C<sub>1</sub>-C<sub>3</sub> alkyl-R<sub>11</sub>)<sub>2</sub>, -CH(CH<sub>2</sub>OH)<sub>2</sub>, -(C<sub>1</sub>-C<sub>3</sub> alkyl)(CH<sub>2</sub>OH)<sub>2</sub>, -(C<sub>1</sub>-C<sub>3</sub> alkyl)-O-(C<sub>1</sub>-C<sub>3</sub> alkyl)-R<sub>11</sub>, -(C<sub>1</sub>-C<sub>3</sub> alkyl)-O-(C<sub>1</sub>-C<sub>3</sub> alkyl)-NH<sub>2</sub>, -(C<sub>1</sub>-C<sub>3</sub> alkyl)-O-(C<sub>1</sub>-C<sub>3</sub> alkyl)-NH-(C<sub>1</sub>-C<sub>3</sub> alkyl)-R<sub>11</sub>, -(C<sub>1</sub>-C<sub>3</sub> alkyl)-O-(C<sub>1</sub>-C<sub>3</sub> alkyl)-N(C<sub>1</sub>-C<sub>3</sub> alkyl-R<sub>11</sub>)<sub>2</sub>, phenyl substituted by one or two groups selected from NH<sub>2</sub>, -N(C<sub>1</sub>-C<sub>3</sub> alkyl), -N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, CN or -(C<sub>1</sub>-C<sub>3</sub> alkyl)-tetrazole, or C<sub>1</sub>-C<sub>6</sub> alkyl,





with each of the alkyl chains of any group in this  $R_4$  definition being optionally substituted by from 1 to 4 OH groups;

$R_7$  in each instance is independently selected from H,  $-NH_2$ ,  $NH(C_1-C_3 \text{ alkyl})$ ,  $N(C_1-C_3 \text{ alkyl})_2$ , or  $C_1-C_3 \text{ alkyl}$ ;

$R_8$  is H, OH or  $C_1-C_3 \text{ alkyl}$ ;

$R_9$  is H, OH,  $-NH_2$ ,  $NH(C_1-C_3 \text{ alkyl})$ , or  $N(C_1-C_3 \text{ alkyl})_2$ ;

$R_{10}$  is H or  $C_1-C_3 \text{ alkyl}$ ;

$R_{11}$  is H, CN, OH,  $NH_2$ , F, or  $CF_3$ ,

or a pharmaceutically acceptable salt or ester form thereof.

2. (Currently amended) A compound of Claim 1 selected from the group of:

1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3-hydroxy-5-methoxy phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(S,S)-1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-[(5-hydroxymethyl-2-phenyl-[1,3]dioxolan-4-ylmethyl)-amino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(S,S)-1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3,4-tri-hydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(4-hydroxy-cyclohexyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2-hydroxy-1-hydroxy-methyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

7-(3-Amino-2-hydroxy-propylamino)-1-cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-[2-(2-hydroxy-ethoxy)-ethylamino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopropyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

7-(4-Diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

or a pharmaceutically acceptable salt or ester form thereof.

3. (Currently amended) A compound of Claim 1 selected from the group of:

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-(2-hydroxy-1-hydroxymethyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-(4-hydroxy-cyclohexylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxyphenyl)-7-(2,3-dihydroxybutylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxypropyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

7-(4-Amino-2,3-dihydroxy-butylamino)-1-cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-([S,S]-2,3,4-triHydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-[2-(2-hydroxyethoxy)-ethyl-amino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

7-(4-Amino-2,3-dihydroxy-butylamino)-3-(2,6-difluoro-3,5-dimethoxyphenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

7-{3-[Bis-(2-hydroxy-ethyl)-amino]-propylamino}-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxypropyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

or a pharmaceutically acceptable salt or ester form thereof.

4. (Currently amended) A compound of Claim 1 selected from the group of:

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxy-butylamino)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

Ethyl-4-[3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-pyrimido[4,5-d]pyrimidin-1-yl]-cyclohexanecarboxylate;

4-[3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-

pyrimido[4,5-d]pyrimidin-1-yl]-cyclohexanecarboxylic acid;

7-Amino-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-piperidin-4-yl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; compound with trifluoroacetic acid; or

1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(2-hydroxy-1-hydroxymethyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Ethyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(4-hydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(S)-1-Cyclopentyl-7-(2,3-dihydroxypropylamino)-3-(2-fluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Ethyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(4-hydroxy-cyclohexylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

4-[3-(2,6-Difluoro-3,5-dimethoxyphenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-pyrimido[4,5-d]pyrimidin-1-yl]-piperidine-1-carboxylic acid; or

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-methylamino-1-piperidin-4-yl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxy-propylamino)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(3-ethoxy-2,6-difluoro-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(R)-1-Cyclopentyl-7-(2,3-dihydroxy-propylamino)-3-(2-fluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Ethyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-[2-(2-hydroxy-ethoxy)-ethylamino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-(*trans*-4-hydroxycyclohexyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

or a pharmaceutically acceptable salt or ester form thereof.

5. A pharmaceutical composition comprising a pharmaceutically effective amount of a compound of Claim 1 and a pharmaceutically acceptable carrier.

6-9. (Canceled)